

Spin Localization of a Fermi Polaron in a Quasirandom Optical Lattice

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Recently, the topics of many-body localization (MBL) and one-dimensional strongly interacting few-body systems have received a lot of interest. These two topics have been largely developed separately. However, the generality of the latter as far as external potentials are concerned – including random and quasirandom potentials – and their shared spatial dimensionality, makes it an interesting way of dealing with MBL in the strongly interacting regime. Utilising tools developed for few-body systems we look to gain insight into the localization properties of the spin in a Fermi gas with strong interactions. We observe a delocalized–localized transition over a range of fillings of a quasirandom lattice. We find this transition to be of a different nature for low and high fillings, due to the diluteness of the system for low fillings.

I. INTRODUCTION

Strongly interacting one-dimensional quantum systems have attracted major attention in recent years [1–5]. When confined to one dimension the fermionic system exhibits a spin-charge separation, and for very strong interactions the charge degrees of freedom are frozen, making it possible to write an effective spin chain Hamiltonian for the system [2–4]. Methods have been developed to solve numerically for the exchange coefficients of this spin chain for any given confining potential [3; 6].

The presence of disorder in an interacting system can result in the violation of the eigenstate thermalization hypothesis, due to many-body localization (MBL) [7; 8]. The localization of single-particle states in the presence of disorder in quantum systems was originally considered by Anderson in 1958 [9]. Over the intervening decades, Anderson localization has been observed in many systems, including in electron gases [10], photonic lattices [11], and cold atoms [12]. For a MBL phase in the tight-binding approximation all eigenstates of the system are Anderson localized [7]. Theoretical work on MBL has been focused on the nature of the delocalization-localization phase transition as disorder is increased [13–16]. Quantum spin chains have been fruitful models for looking at this transition. In most cases the disorder is introduced in the external magnetic field or coupling coefficients of the spin chain. In this work we still consider a quantum spin chain, but one that is induced by the strong interactions present between fermions. We introduce disorder in the system via a quasirandom optical lattice potential.

In recent years, the field of ultracold atomic gases in one dimension has received a lot of interest [17–19]. Such systems have been considered for strongly interacting fermions [20] and bosons [21–23]. In this field, the MBL phase transition has been observed with interacting fermions in a one dimensional quasirandom optical lattice [24]. Recently, an experimental realization of only a few strongly interacting fermions in a one dimensional

trap has been realised [25].

II. MODEL

We consider N strongly interacting repulsive fermions in one dimension (see Fig. 1d). This system is described by the Hamiltonian,

$$H = \sum_{i=1}^N \left(\frac{p_i^2}{2m} + V(x_i) \right) + g \sum_{i < j}^N \delta(x_i - x_j) \quad (1)$$

where g is the interaction strength, and $V(x_i)$ is the single-particle external potential. We consider the repulsive case $g > 0$, and take $g \rightarrow +\infty$.

As discussed above, strongly interacting spin-1/2 fermions in one dimension can effectively be written as a spin chain Hamiltonian. The coefficients of the spin chain have been found to be dependent on the potential and the single particle wavefunctions [2–4; 6], which are solutions to the first term in the Hamiltonian (Eq. (1)). We use the open source program CONAN [6], which has been developed to take arbitrary potentials and numerically calculate the $N - 1$ exchange coefficients between the spin chain sites for up to $N \sim 30$ particles.

We consider a quasirandom potential with open boundary conditions, with a main lattice of strength V_1 and a disorder term of strength V_2 . The potential $V(x)$, appearing in Eq. (1) is given by

$$V(x) = V_1 \cos(\tau_1 L_s x) + V_2 \cos(\tau_2 L_s x + \phi), \quad (2)$$

where L_s is the number of wells – or ‘sites’ – in the lattice, and x is defined in units of L_s . Throughout this work we set $\tau_1 = 2\pi$ and $\tau_2 = 1$, satisfying the need for τ_1/τ_2 to be incommensurate for the above potential to be quasirandom. We will quantify the disorder strength by the ratio $W = V_2/V_1$. We calculate the spin chain coefficients for several realizations of each W with various phases, ϕ ,

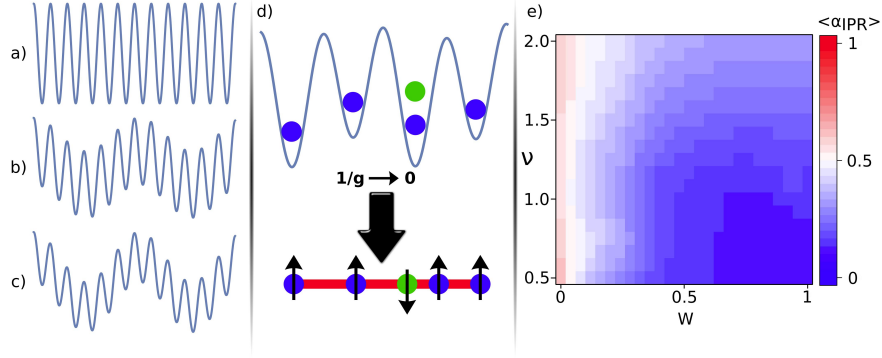


FIG. 1. a – c) The quasirandom potential, Eq. (2), for $W = 0, 0.5$ and 1 respectively. d) Illustration of the mapping to an effective spin chain model for strong interactions. e) Average inverse participation ratio for a filling $\nu = N/L_s$ of the N single particle states for disorder strength W .

then average over the realizations for the final observables. We consider the disorder range of $0 \leq W \leq 1$, with examples shown in Fig. 1a, b and c. The main lattice strength V_1 is chosen to ensure that the lattice is strong enough to be felt by all particles, without the particles being localized into single sites. Defining the unit of energy as $1/(2L_s^2)$, having put $\hbar = m = 1$, we set the lattice strength to the dimensionless quantity $V_1 = 10L_s^2$ for all calculations.

For the spin chain we have the Hamiltonian

$$K = -\frac{1}{2} \sum_{j=1}^{N-1} J_j (\boldsymbol{\sigma}_j \cdot \boldsymbol{\sigma}_{j+1} - 1), \quad (3)$$

where $\boldsymbol{\sigma}_j$ are the Pauli matrices acting on the j th site, and J_j is the coefficient connecting the j and $j+1$ sites. The spin chain has hard wall boundaries, see Fig. 1d. The energy of the physical system is given in the groundstate manifold by

$$E_n = E_0 - \frac{K_n}{g}, \quad (4)$$

with E_0 being the many-body groundstate energy for $g = \infty$, i.e. the sum of the first N single particle eigenenergies, and K_n the eigenvalues of the spin chain Hamiltonian. As we will consider the ordering of the states, and the relative energy between them, we only require knowledge of the spin chain eigenvalues $-K_n$. As already stated, we use the program CONAN [6] to compute the spin chain coefficients J_j over a range of W , for a quasirandom potential with $L_s = 12$ lattice sites, and a range of particle numbers $N = 6, 7, 8, \dots, 24$, corresponding to fillings $\nu = 1/2, 0.583, 0.667, \dots, 2$, with the definition $\nu = N/L_s$. For each W and N we average over 19 realizations of various phases ϕ . Using the calculated spin chain coefficients, we solve the stationary Schrödinger equation for the spin chain Hamiltonian, Eq. (3), numerically. For

the polaron we will denote the wavefunction as

$$|\Psi\rangle = \sum_{j=1}^N C_j |\uparrow \dots \uparrow (\downarrow)_j \uparrow \dots \uparrow\rangle, \quad (5)$$

where C_j is the coefficient for the polaron in the j th spin chain site. To gain further insight, we will also consider the case of two polarons, which we expect to have similar general behaviour to the single polaron in this system. For two polarons we write the wavefunction as

$$|\Psi\rangle = \sum_{i < j}^N C_{(i,j)} |\uparrow \dots \uparrow (\downarrow)_i \uparrow \dots \uparrow (\downarrow)_j \uparrow \dots \uparrow\rangle. \quad (6)$$

III. MEASURES OF LOCALIZATION

The onset of Anderson localization in the system can be observed by considering the inverse participation ratio (IPR) [5], given by

$$\alpha_{\text{IPR}} = \frac{1}{m \sum_v |C_v|^4}, \quad (7)$$

with C_v being the coefficients of either the polaron $v = j$ or two polaron $v = (i, j)$ states, and m denoting the size of the Hilbert space of the wavefunction. For a fully delocalized state $\alpha_{\text{IPR}} \sim 1$. For a fully localized state we have a convergence of α_{IPR} towards zero. We will consider the IPR of the ground, and the highest energy states. In addition we calculate the average IPR across all other states (denoted by $\langle \alpha_{\text{IPR}} \rangle$). The average IPR gives an indication of the overall localization of the system.

A standard measure of the MBL transition is the properties of the energy level statistics, which can be investigated via the ratio of adjacent energy level gaps [5; 15; 26; 27]

$$r_n = \frac{\min(\delta_n, \delta_{n-1})}{\max(\delta_n, \delta_{n-1})}, \quad (8)$$

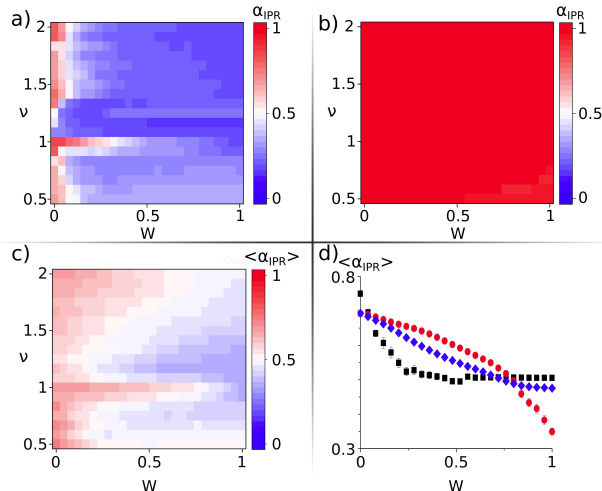


FIG. 2. α_{IPR} for the single polaron spin chain. a) Groundstate IPR. b) highest energy state IPR. c) The average IPR over all states except the ground and highest, $\langle\alpha_{\text{IPR}}\rangle$. d) cut-outs of c) for $\nu = 0.5$ squares (black), $\nu = 1$ circles (red), and $\nu = 2$ diamonds (blue).

where $\delta_n = E_n - E_{n-1}$ is the gap in the spectrum between the E_n and E_{n-1} eigenvalues, with min and max taking the minimum and maximum value of the two adjacent gaps in the spectrum (δ_n, δ_{n-1}), ensuring $0 \leq r_n \leq 1$. We will take the average of the gap ratio $\langle r_n \rangle$ over all δ_n , with the exclusion of the ground and highest energy states as will be discussed in Sec. IV. In the delocalized phase we expect the energy level statistics to satisfy a Wigner-Dyson distribution (WD) [15; 26], with the average ratio $\langle r_n \rangle_{\text{WD}} \simeq 0.536$ [5]. Meanwhile, the MBL phase has statistics that satisfy the Poisson distribution (PD) [15; 26], with an average ratio of $\langle r_n \rangle_{\text{PD}} \simeq 0.386$ [5].

IV. LOCALIZATION OF THE SPIN

First we confirm the localization in the charge degree of freedom. In Fig. 1e, we consider the average IPR over all single particle states for the quasirandom lattice, Eq. (2). We find for disorder above $W \sim 0.5$ the general localization of the single particle states across the whole range of ν . The critical disorder for the delocalized to localized transition of the single particle states is weakly dependent on the filling of the lattice. This result gives a good indication that the particles are “feeling” the lattice potential.

For the spin of a single polaron we consider α_{IPR} of the ground, and highest energy states, then $\langle\alpha_{\text{IPR}}\rangle$ over all other states, with the results shown in Fig. 2. In addition, we consider two polarons in the system, with the same observables as for the single polaron, but in addition we calculate the average energy gap ratio $\langle r_n \rangle$.

We do not consider the gap ratio for the single polaron due to the small number of states, N , in the system, resulting in large variances in $\langle r_n \rangle$ over the realizations of the disorder. Naturally, for two polarons there is a larger number of states, $N(N-1)/2$, resulting in smaller variance over the disorder realizations.

The groundstate of the spin is found to localize at small disorder in Figs. 2a and 3a, with strong localization for $W > 0.1$, for most ν . With the exception of around unit filling, where we have a spin in each lattice site, resulting in an elongated transition to the localized state. The highest energy state is delocalized across the system over all disorder, Figs. 2b and 3b. Therefore, for our system we can never have a true MBL phase, in the sense that all states will not localize. However, with the inherently delocalized highest energy state excluded, we observe a delocalized-localized transition over a range of ν .

With the average IPR over all states in the system except the ground and highest states, $\langle\alpha_{\text{IPR}}\rangle$, we can gain an insight into the general localization properties of the system, see Figs. 2c and 3c. We observe a defined transition from a majority of states being delocalized to heavily localized over a range of fillings from $1 \leq \nu \leq 2$. For small fillings, $\nu < 1$, we observe a trend towards localization with large disorder. The relatively weak localization of states at these fillings is due to the diluteness of the system. Each fermion (or groups of fermions) can be well separated from its neighbours, resulting in weak coupling coefficients, effectively resulting in the separation of the spin chain into sections. Hence we observe some localization of the state, but not due to disorder in the spin chain. The regimes discussed are well shown by the cut outs of Figs. 2d and 3e.

However, the IPR is a poor measure of the localization of all states, and a standard measure for this (the MBL phase) is the average energy gap ratio in the spectrum, $\langle r_n \rangle$. We calculated $\langle r_n \rangle$ for two polarons in the system, Fig. 3d, with a cut out at select fillings in Fig. 3f. For $\nu \sim 1$, we see a transition from an extended ($\langle r_n \rangle_{\text{WD}} \simeq 0.536$) to a localized phase ($\langle r_n \rangle_{\text{PD}} \simeq 0.386$). At $\nu = 1.0833 = 1 + 1/12$, where we are at one particle over unit filling, we observe the states to have Poisson statistics independent of disorder, shown by the white region above unit filling in Fig. 3d. This is due to the spin chain coefficients having a form that is ‘well-like’ at this filling without the presence of disorder [28]. Thus the statistics of the eigenvalue gaps are that of the Poisson distribution, as has been shown for interacting trapped bosons in harmonic potentials [29]. For higher filling, we see a transition from a delocalized to a localized phase with increasing disorder. However as we approach $\nu = 2$, $\langle r_n \rangle$ is consistently at an intermediate value, Fig. 3f, which is consistent with a mixed phase of localized and delocalized states.

With $\nu < 1$, we observe a different regime of the system. $\langle r_n \rangle$ converges to a value well below 0.386, as seen in Figs. 3d and f, with a weak localization across all states as seen in Figs. 2c and 3c. The convergence value of $\langle r_n \rangle$ is not consistent with any spectrum we know of. For

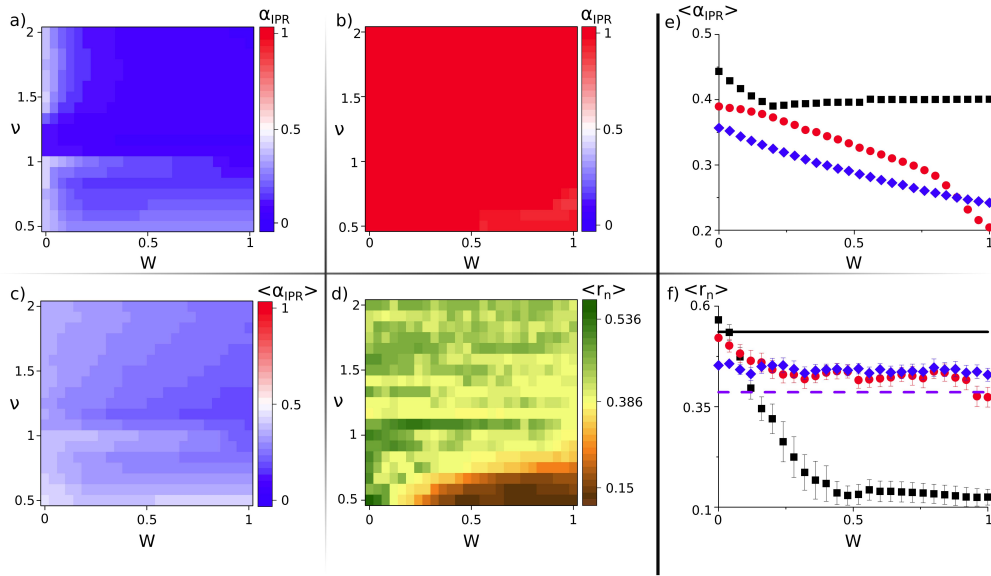


FIG. 3. α_{IPR} for the two-polaron spin chain. a) Groundstate IPR, averaged over 9 disorder phase, ϕ , realizations. b) Same as a) but for the highest energy state. c) The average IPR over all states except the ground and highest, $\langle \alpha_{\text{IPR}} \rangle$. d) The energy gap ratio, $\langle r_n \rangle$. e) and f) cut-outs of c) and d) respectively for $\nu = 0.5$ squares (black), $\nu = 1$ circles (red), and $\nu = 2$ diamonds (blue). On plot f) black solid line and purple dashed line denote the expected values for an extended and localized phase respectively.

$\nu < 1$ the states are localized because of the break up of the spin chain due to the diluteness of the system, and not through disorder. The gap ratio further reflects the different nature of the localization transition of the states for low filling.

V. CONCLUSIONS

Using recent advances in describing strongly interacting confined particles in one dimension, we have investigated the localization properties of the spin degree of freedom. It is well known that the charge degree of freedom is localized in this system in the presence of strong interactions. For the spin we observe the localization of the majority of states upon sufficient disorder for $\nu > 1$. For small fillings, $\nu < 1$, we observe a weak localization regime due to the system being dilute. The system considered can never be completely localized, due to the

presence of a fully delocalized highest energy state. This state is an inherent property of the system. However with the exclusion of this delocalized state, we observe a delocalized to localized transition for both the polaron and two polaron systems. This transition is seen for fillings above unity by the convergence of the level statistics to the Poisson distribution expected in the MBL phase. For low fillings and above a certain disorder strength we see the emergence of a regime with different statistics, due to the diluteness of the system.

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